

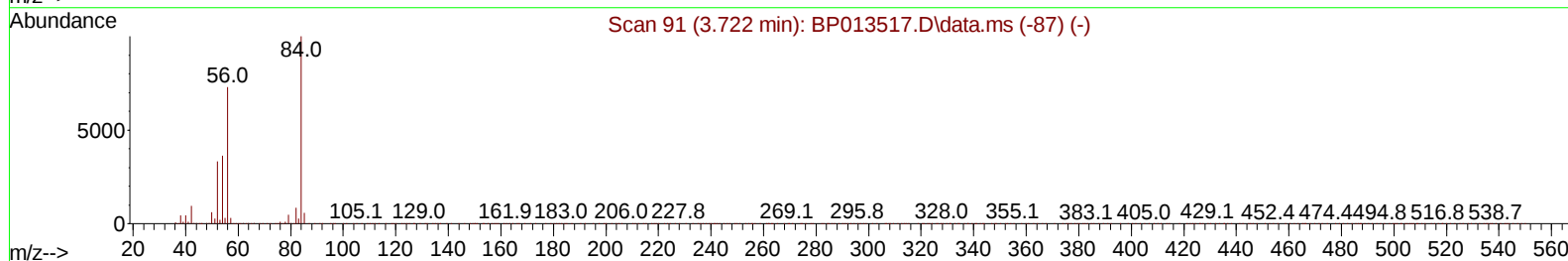
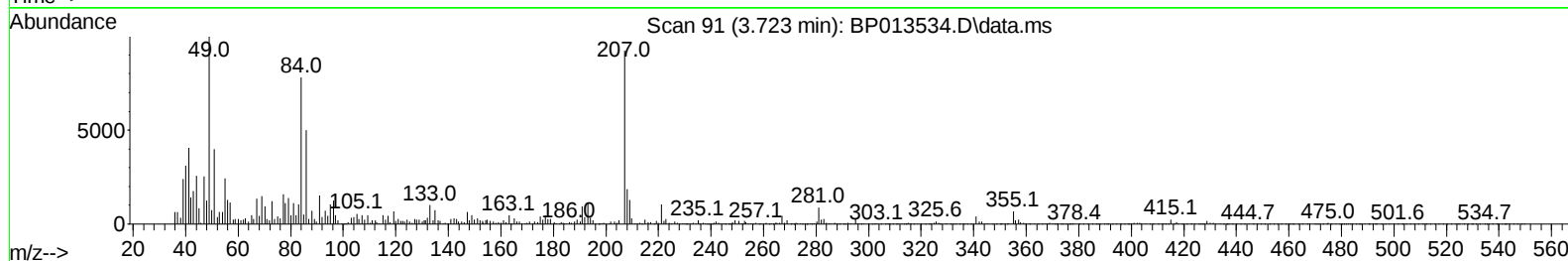
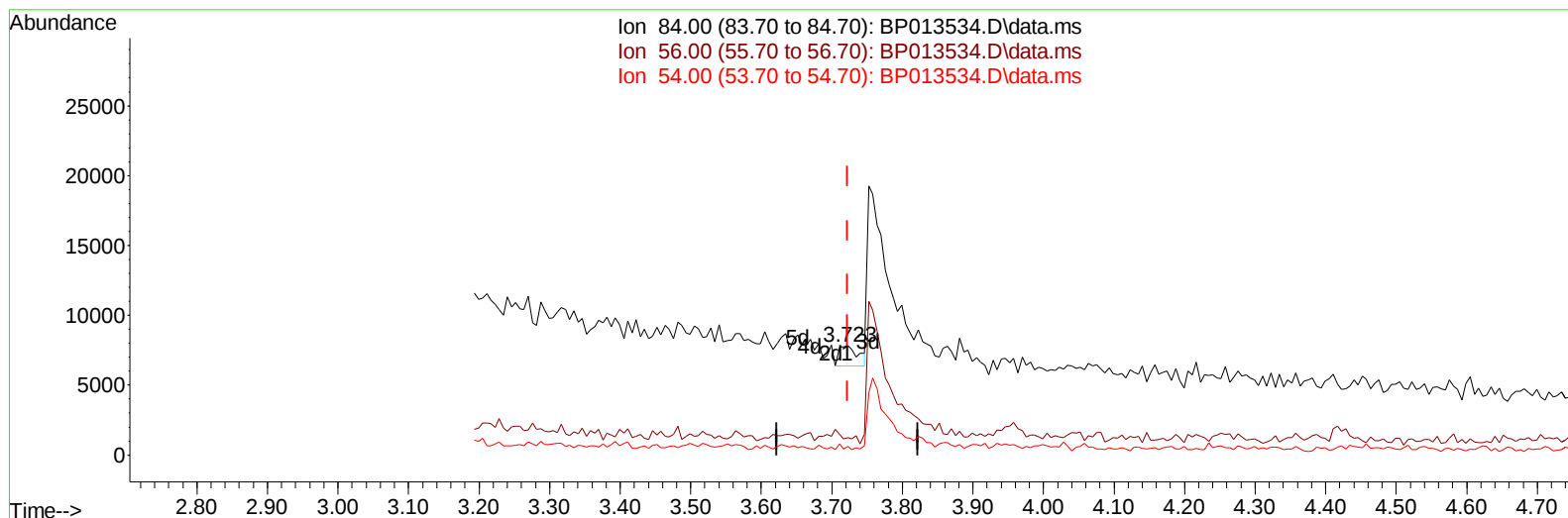
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011823\
 Data File : BP013534.D
 Acq On : 18 Jan 2023 22:12
 Operator : CG/JU
 Sample : 01146-12
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4418

Manual IntegrationsAPPROVED

Quant Time: Jan 19 01:49:39 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 18 01:06:35 2023
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/19/2023
 Supervised By :mohammad ahmed 01/20/2023



TIC: BP013534.D\data.ms

(4) Pyridine-d5 (S)

3.723min (+ 0.000) 0.06 ng/u1

response 2694

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	72.30	16.22#
54.00	34.30	8.09#
0.00	0.00	0.00

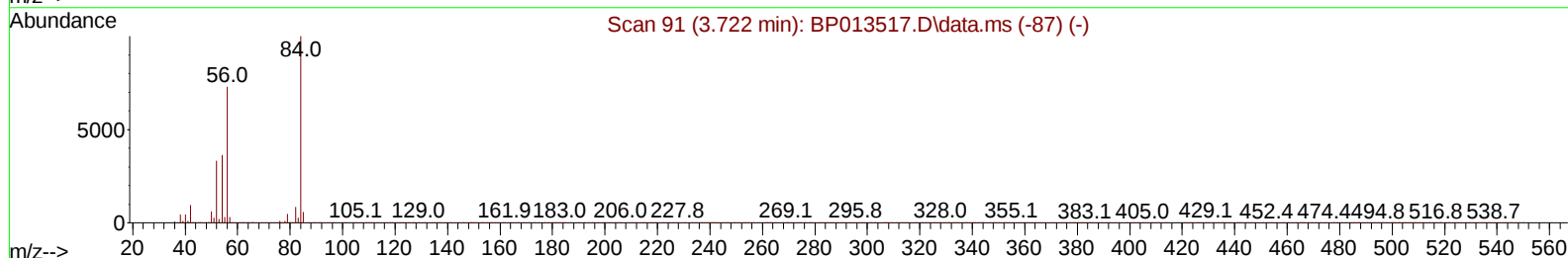
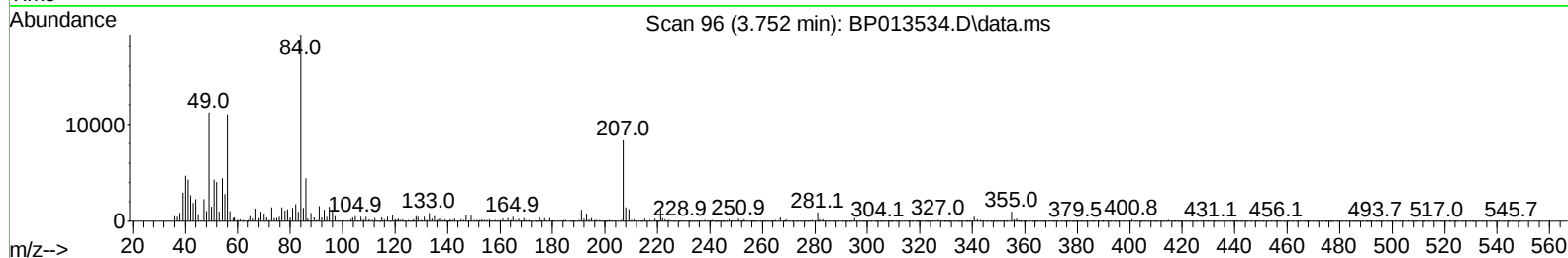
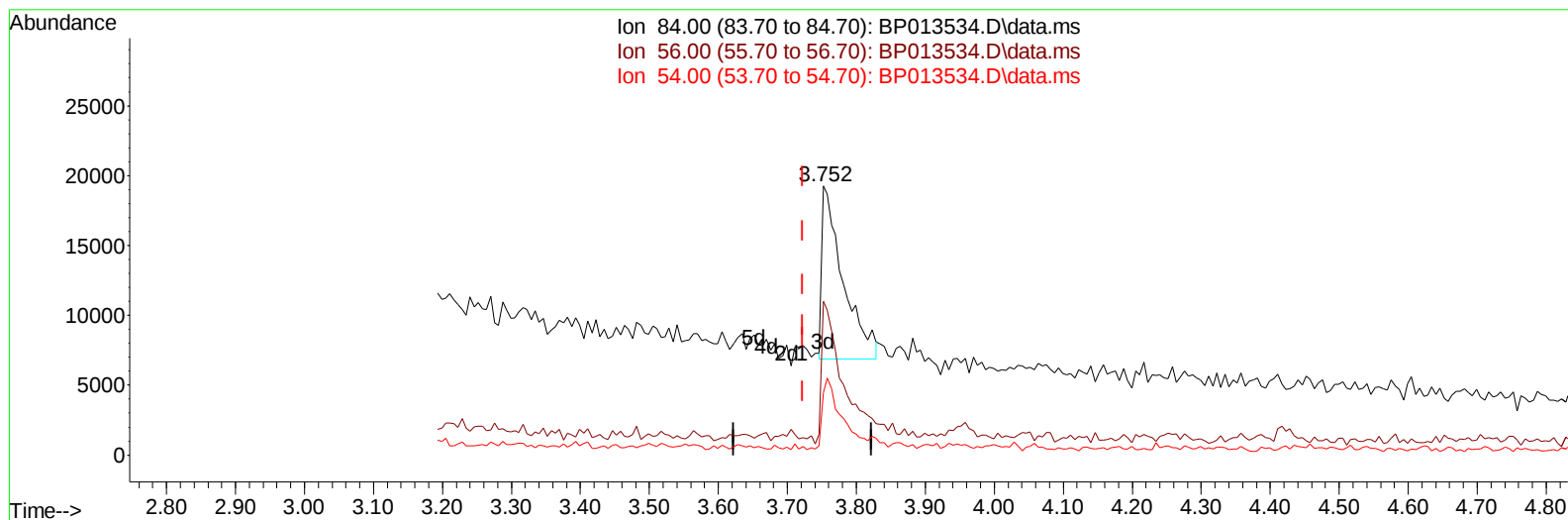
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TIC: BP013534.D\data.ms

(4) Pyridine-d5 (S)

3.752min (+ 0.030) 0.56 ng/u1 m

response 26402

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	72.30	57.18#
54.00	34.30	23.16#
0.00	0.00	0.00

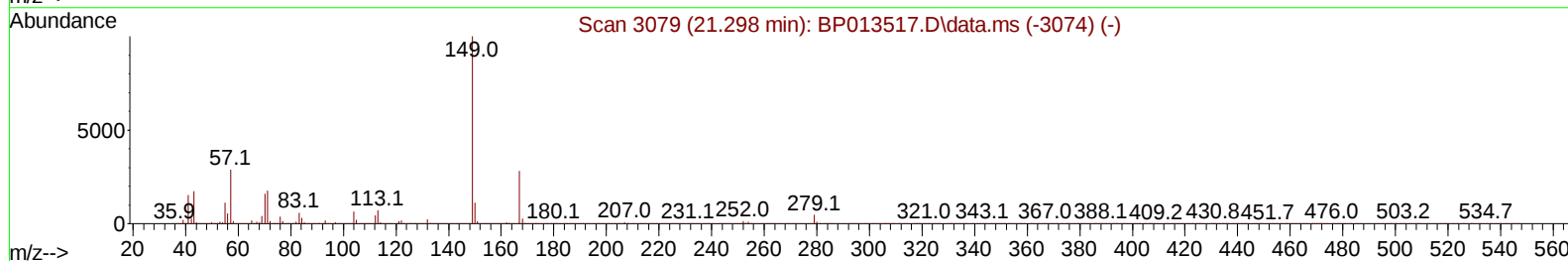
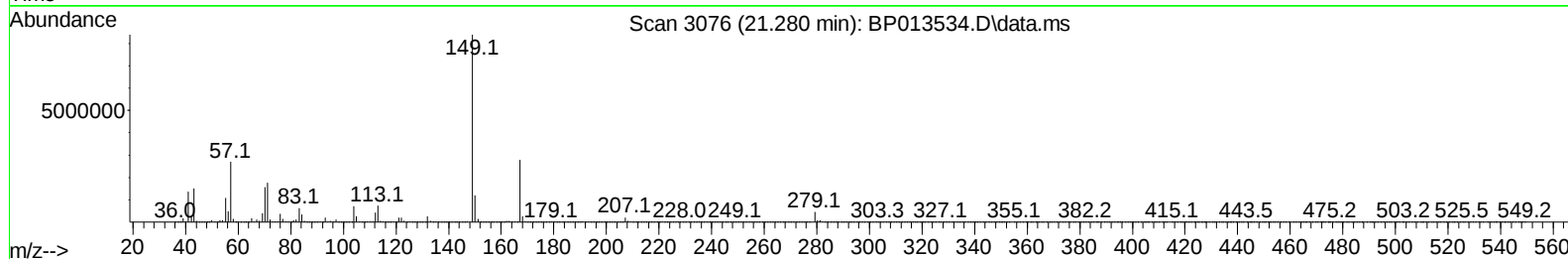
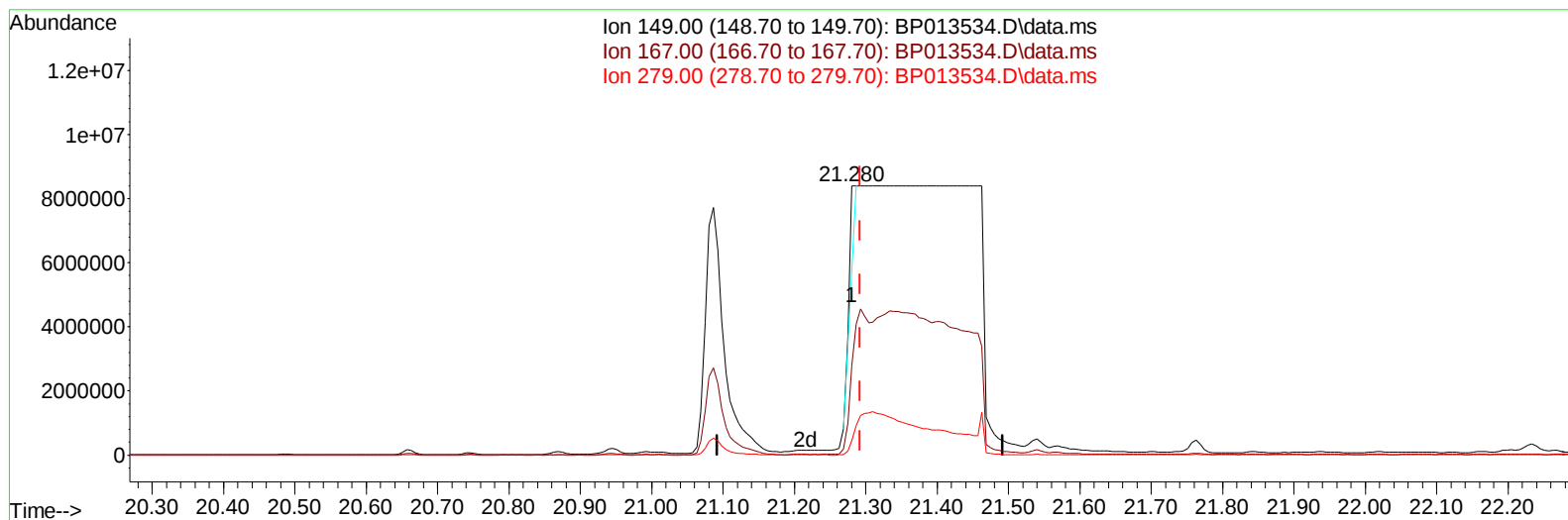
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 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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TIC: BP013534.D\data.ms

(86) Bis(2-ethylhexyl)phthalate

21.280min (-0.012) 1345.85 ng/u1

response 2595222

Ion	Exp%	Act%
149.00	100.00	100.00
167.00	27.80	33.21
279.00	4.40	5.39#
0.00	0.00	0.00

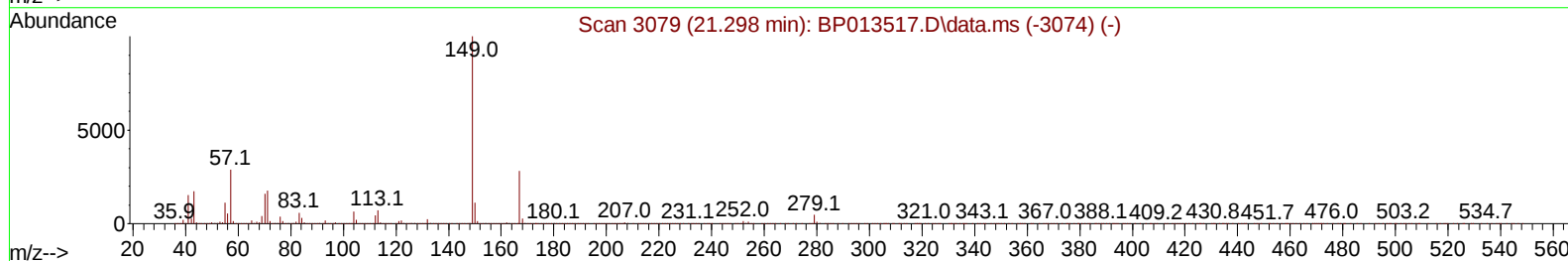
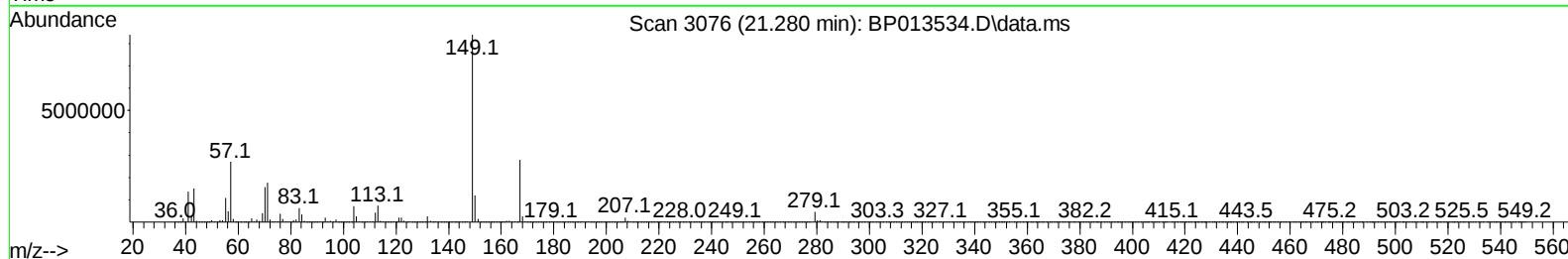
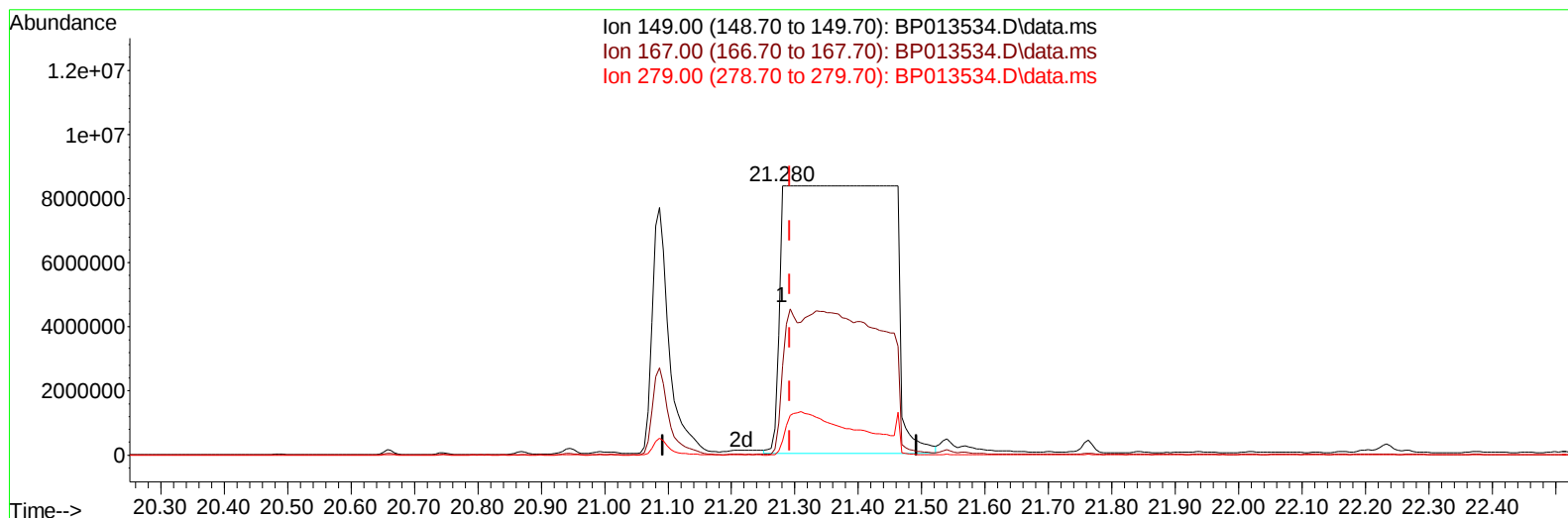
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TIC: BP013534.D\data.ms

(86) Bis(2-ethylhexyl)phthalate

21.280min (-0.012) 50547.13 ng/ul m

response 97470596

Ion	Exp%	Act%
149.00	100.00	100.00
167.00	27.80	33.21
279.00	4.40	5.39#
0.00	0.00	0.00

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Instrument :
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.899	152		652694	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136		2442078	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.546	164		1280833	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188		2952774	20.000	ng/ul	0.00
79) Chrysene-d12	21.392	240		69141	20.000	ng/ul	# 0.00
88) Perylene-d12	23.857	264		2885004	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.299	96		40288	2.399	ng/uL	0.00
4) Pyridine-d5	3.752	84		26402m	0.564	ng/ul	0.03
7) Phenol-d5	7.058	99		981533	18.123	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67		572719	16.762	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132		776591	19.227	ng/ul	0.00
15) 4-Methylphenol-d8	8.611	113		774281	18.765	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128		358956	20.644	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143		421629	22.151	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165		761637	20.078	ng/ul	0.00
31) 4-Chloroaniline-d4	10.852	131		465844	8.717	ng/ul	0.00
46) Dimethylphthalate-d6	13.951	166		1900277	20.364	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160		2197294	21.272	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143		214040	12.531	ng/ul	0.00
60) Fluorene-d10	15.540	176		1539214	18.674	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200		211802	11.312	ng/ul	0.02
73) Anthracene-d10	17.410	188		2184887	16.741	ng/ul	0.00
81) Pyrene-d10	19.639	212		2723095	723.876	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264		2963791	20.914	ng/ul	0.00
Target Compounds							
86) Bis(2-ethylhexyl)phtha...	21.280	149		97470596m	50547.125	ng/ul	Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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